



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:21 AM UTC

PDB ID : 9UDE / pdb_00009ude
Title : Crystal structure of MonCI in complex with diepoxidized farnesyl acetate
Authors : Xiao, H.L.; Guan, Y.Z.; Chen, X.
Deposited on : 2025-04-06
Resolution : 2.65 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

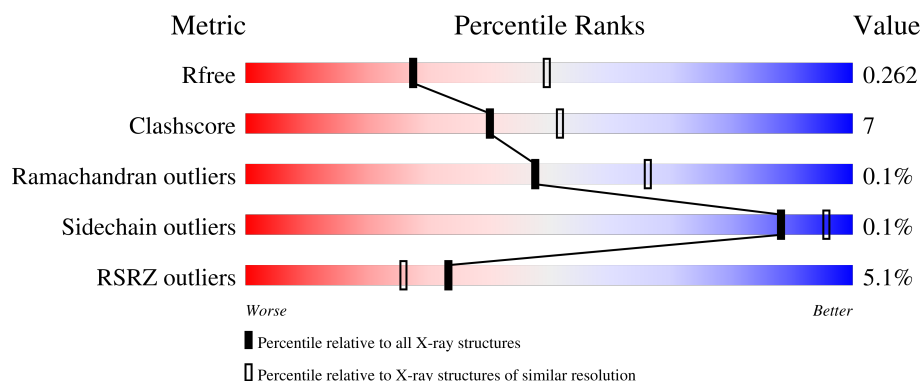
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	511	
1	B	511	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7237 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called MonCI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	472	Total	C	N	O	S	0	1	0
			3547	2219	645	673	10			
1	B	472	Total	C	N	O	S	0	0	0
			3503	2196	638	660	9			

There are 30 discrepancies between the modelled and reference sequences:

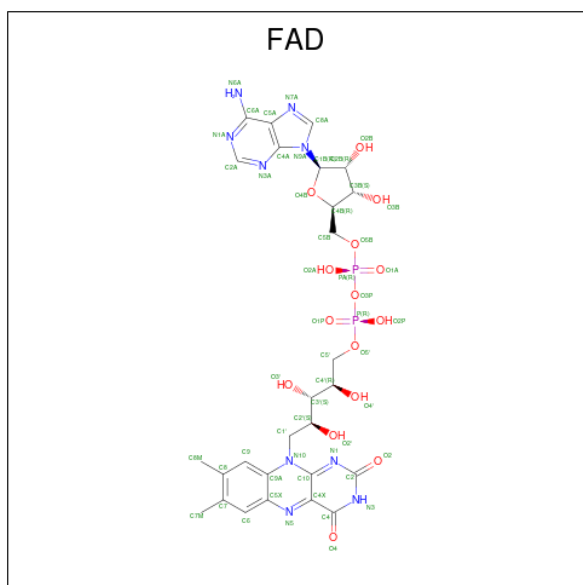
Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	MET	-	initiating methionine	UNP Q846W9
A	-13	ASN	-	expression tag	UNP Q846W9
A	-12	HIS	-	expression tag	UNP Q846W9
A	-11	LYS	-	expression tag	UNP Q846W9
A	-10	VAL	-	expression tag	UNP Q846W9
A	-9	HIS	-	expression tag	UNP Q846W9
A	-8	HIS	-	expression tag	UNP Q846W9
A	-7	HIS	-	expression tag	UNP Q846W9
A	-6	HIS	-	expression tag	UNP Q846W9
A	-5	HIS	-	expression tag	UNP Q846W9
A	-4	HIS	-	expression tag	UNP Q846W9
A	-3	ILE	-	expression tag	UNP Q846W9
A	-2	GLU	-	expression tag	UNP Q846W9
A	-1	GLY	-	expression tag	UNP Q846W9
A	0	ARG	-	expression tag	UNP Q846W9
B	-14	MET	-	initiating methionine	UNP Q846W9
B	-13	ASN	-	expression tag	UNP Q846W9
B	-12	HIS	-	expression tag	UNP Q846W9
B	-11	LYS	-	expression tag	UNP Q846W9
B	-10	VAL	-	expression tag	UNP Q846W9
B	-9	HIS	-	expression tag	UNP Q846W9
B	-8	HIS	-	expression tag	UNP Q846W9
B	-7	HIS	-	expression tag	UNP Q846W9
B	-6	HIS	-	expression tag	UNP Q846W9
B	-5	HIS	-	expression tag	UNP Q846W9

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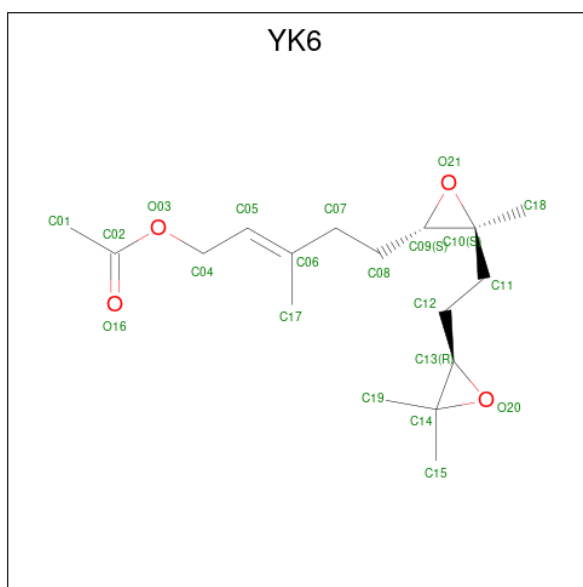
Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	HIS	-	expression tag	UNP Q846W9
B	-3	ILE	-	expression tag	UNP Q846W9
B	-2	GLU	-	expression tag	UNP Q846W9
B	-1	GLY	-	expression tag	UNP Q846W9
B	0	ARG	-	expression tag	UNP Q846W9

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$) (labeled as "Ligand of Interest" by depositor).



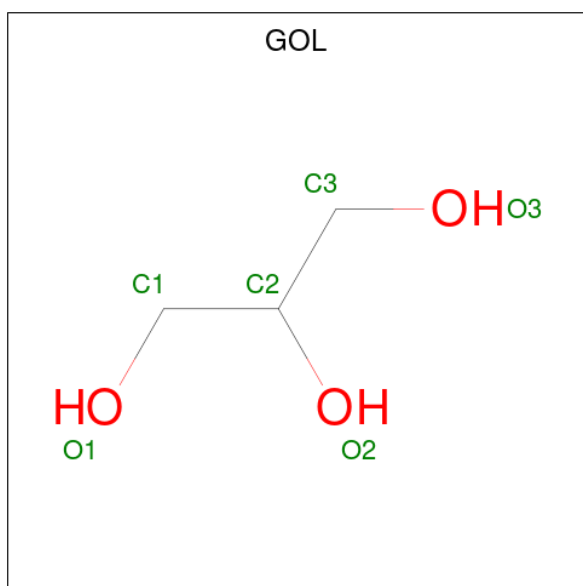
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is diepoxidized farnesyl acetate (CCD ID: YK6) (formula: $C_{17}H_{28}O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			21	17	4		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	Cl 1	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	45	Total 45	O 45	0	0
6	B	8	Total 8	O 8	0	0

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- Molecule 1: MonCl



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.62Å 52.77Å 145.83Å 90.00° 92.37° 90.00°	Depositor
Resolution (Å)	145.70 – 2.65 145.70 – 2.65	Depositor EDS
% Data completeness (in resolution range)	95.7 (145.70-2.65) 95.7 (145.70-2.65)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 2.65Å)	Xtriage
Refinement program	PHENIX (1.16_3549: ???)	Depositor
R, R_{free}	0.199 , 0.261 0.203 , 0.262	Depositor DCC
R_{free} test set	1423 reflections (4.69%)	wwPDB-VP
Wilson B-factor (Å ²)	46.5	Xtriage
Anisotropy	0.491	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 53.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.025 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7237	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.63% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FAD, CL, YK6, GOL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.17	0/3620	0.41	0/4939
1	B	0.20	0/3573	0.49	0/4881
All	All	0.18	0/7193	0.45	0/9820

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3547	0	3502	36	0
1	B	3503	0	3443	58	0
2	A	53	0	31	2	0
2	B	53	0	31	1	0
3	A	21	0	0	0	0
4	A	6	0	8	0	0
5	A	1	0	0	1	0
6	A	45	0	0	0	0
6	B	8	0	0	0	0
All	All	7237	0	7015	94	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including

hydrogen atoms). The all-atom clashscore for this structure is 7.

All (94) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:334:HIS:CE1	1:B:382:LEU:HD23	1.85	1.10
1:B:334:HIS:CE1	1:B:382:LEU:CD2	2.56	0.88
1:A:233:ARG:HH21	1:A:411:LEU:HD23	1.39	0.87
1:B:195:PRO:HG2	1:B:311:LEU:HD11	1.61	0.81
1:B:334:HIS:HE1	1:B:382:LEU:HD23	1.47	0.80
1:B:176:ILE:HD13	1:B:319:LEU:HB2	1.68	0.74
1:A:469:PRO:HG2	1:A:470:GLU:OE2	1.88	0.73
1:B:4:THR:HG22	1:B:5:ARG:H	1.56	0.71
1:B:36:GLU:HG2	1:B:40:LEU:HD21	1.73	0.70
1:B:451:MET:HE3	1:B:455:ALA:HB2	1.76	0.68
1:B:211:ARG:HG3	1:B:295:VAL:HG12	1.79	0.64
1:B:212:LEU:HD21	1:B:250:ARG:HD2	1.79	0.64
1:B:141:VAL:O	1:B:188:TRP:NE1	2.32	0.62
1:A:407:THR:O	1:A:411:LEU:HD12	2.00	0.61
1:B:189:LEU:HD22	1:B:194:VAL:HG21	1.82	0.59
1:B:309:GLU:HB2	1:B:370:LYS:HG2	1.85	0.59
1:B:230:ALA:H	1:B:240:PHE:HD1	1.52	0.58
1:A:4:THR:OG1	1:A:5:ARG:N	2.35	0.56
1:B:38:ASP:HA	1:B:137:ARG:NH2	2.21	0.55
1:A:233:ARG:HG2	1:A:390:TYR:CE1	2.43	0.54
1:A:208:TYR:CZ	1:A:330:PRO:HG2	2.43	0.53
1:B:334:HIS:HE1	1:B:382:LEU:CD2	2.12	0.53
1:A:468:LEU:HD23	1:A:470:GLU:OE1	2.08	0.53
1:A:23:HIS:HB2	1:A:126:ALA:HB1	1.92	0.52
1:A:37:ARG:HD3	1:A:139:GLU:HG3	1.92	0.52
1:B:202:VAL:HG22	1:B:395:VAL:HG23	1.90	0.52
1:B:208:TYR:CE1	1:B:330:PRO:HG2	2.44	0.52
1:B:353:SER:C	1:B:355:GLN:H	2.18	0.52
1:B:381:ILE:HD13	1:B:401:ARG:CZ	2.39	0.52
1:B:32:VAL:O	1:B:132:ILE:HA	2.10	0.51
1:B:95:LEU:HB2	1:B:100:TRP:CZ3	2.46	0.50
1:A:5:ARG:HH21	1:A:5:ARG:HG2	1.77	0.50
1:A:233:ARG:HG2	1:A:390:TYR:CZ	2.46	0.50
1:B:38:ASP:OD2	1:B:49:GLY:N	2.45	0.50
1:B:179:THR:OG1	1:B:183:SER:HB2	2.11	0.50
1:A:208:TYR:OH	1:A:331:ILE:HG13	2.12	0.49
1:B:206:VAL:HB	1:B:301:GLY:HA3	1.94	0.49
1:B:233:ARG:HG3	1:B:390:TYR:CE1	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:208:TYR:CZ	1:B:330:PRO:HG2	2.47	0.49
1:B:259:ARG:NH1	1:B:392:ASN:O	2.44	0.49
1:B:305:ARG:HH11	1:B:380:TRP:CD1	2.31	0.49
1:B:472:LEU:HD22	1:B:477:LEU:HB2	1.95	0.49
1:B:10:VAL:HG23	1:B:172:ALA:HB2	1.95	0.48
1:A:242:VAL:HG11	1:A:244:TYR:CE2	2.48	0.48
1:A:470:GLU:H	1:A:470:GLU:CD	2.20	0.48
1:B:177:ASP:OD1	1:B:179:THR:OG1	2.26	0.48
1:A:92:LEU:HG	1:A:225:ALA:HB3	1.95	0.48
1:A:338:SER:HB2	1:A:379:PRO:HG3	1.95	0.48
1:B:188:TRP:O	1:B:192:LEU:HD13	2.14	0.48
1:B:407:THR:O	1:B:411:LEU:HD12	2.14	0.48
1:A:34:VAL:HB	1:A:134:VAL:HG12	1.94	0.47
1:B:4:THR:HG22	1:B:5:ARG:HG3	1.96	0.47
1:B:232:ASP:OD1	1:B:233:ARG:N	2.48	0.47
1:B:258:THR:HG21	1:B:390:TYR:CE1	2.49	0.47
1:B:161:ASP:HB3	1:B:164:SER:HG	1.80	0.47
1:A:243:VAL:HG23	1:A:282:LEU:HD22	1.97	0.46
1:A:208:TYR:CE1	1:A:330:PRO:HG2	2.50	0.46
1:B:38:ASP:HA	1:B:137:ARG:HH21	1.79	0.46
1:A:37:ARG:HG3	1:A:162:LEU:HD11	1.98	0.46
1:B:382:LEU:HD11	1:B:434:VAL:HG12	1.97	0.46
1:B:382:LEU:HD12	1:B:441:GLN:HG3	1.98	0.46
1:A:298:SER:OG	5:A:504:CL:CL	2.69	0.46
1:A:95:LEU:HB2	1:A:100:TRP:CZ3	2.51	0.46
1:A:349:GLU:O	1:A:353:SER:OG	2.25	0.45
1:B:208:TYR:OH	1:B:331:ILE:HG13	2.16	0.45
1:A:319:LEU:HD11	1:A:368:LEU:HD13	1.98	0.45
1:A:258:THR:HG21	1:A:390:TYR:CE1	2.52	0.45
1:B:259:ARG:HD2	1:B:392:ASN:HB3	1.99	0.45
1:B:212:LEU:HD13	1:B:252:LEU:HD13	1.98	0.45
1:B:385:THR:HG21	1:B:441:GLN:HE22	1.82	0.45
1:B:211:ARG:CG	1:B:295:VAL:HG12	2.47	0.44
1:A:447:ASN:HB3	1:A:451:MET:HE3	2.00	0.44
1:B:181:ARG:HG2	1:B:304:ARG:NH2	2.33	0.44
1:B:144:THR:HG21	1:B:169:GLN:HE21	1.82	0.43
1:A:55:HIS:NE2	1:A:298:SER:HB2	2.34	0.43
1:B:161:ASP:HB3	1:B:164:SER:OG	2.18	0.43
1:A:55:HIS:ND1	1:A:210:THR:OG1	2.45	0.43
1:B:189:LEU:HB3	1:B:194:VAL:HG23	2.01	0.43
1:B:269:GLU:C	1:B:272:PRO:HD2	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:VAL:HG21	1:B:244:TYR:HB3	1.99	0.43
1:B:57:HIS:HA	2:B:700:FAD:O4	2.19	0.42
1:B:247:GLU:OE1	1:B:250:ARG:NH2	2.53	0.42
1:A:242:VAL:HG11	1:A:244:TYR:CZ	2.55	0.41
1:B:17:ALA:HB1	1:B:321:ILE:HD12	2.01	0.41
1:A:335:GLY:HA3	2:A:501:FAD:H1'2	2.03	0.41
1:B:243:VAL:HG23	1:B:282:LEU:HD22	2.03	0.41
1:B:92:LEU:HG	1:B:225:ALA:HB3	2.02	0.41
1:B:381:ILE:HD13	1:B:401:ARG:NE	2.36	0.41
1:A:305:ARG:HH11	1:A:380:TRP:CD1	2.39	0.41
1:A:269:GLU:C	1:A:272:PRO:HD2	2.46	0.40
1:A:446:SER:O	1:A:450:LEU:HG	2.21	0.40
1:A:472:LEU:HD13	1:A:478:ASP:HA	2.03	0.40
1:A:206:VAL:HA	1:A:257:CYS:O	2.22	0.40
1:A:57:HIS:HA	2:A:501:FAD:O4	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	469/511 (92%)	454 (97%)	15 (3%)	0	100	100
1	B	468/511 (92%)	452 (97%)	15 (3%)	1 (0%)	43	60
All	All	937/1022 (92%)	906 (97%)	30 (3%)	1 (0%)	48	66

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	354	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	361/401 (90%)	361 (100%)	0	100	100
1	B	351/401 (88%)	350 (100%)	1 (0%)	86	94
All	All	712/802 (89%)	711 (100%)	1 (0%)	88	95

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	232	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	125	GLN
1	B	169	GLN
1	B	438	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	YK6	A	502	-	22,22,22	3.02	5 (22%)	33,33,33	3.73	9 (27%)
2	FAD	B	700	-	58,58,58	0.27	0	85,89,89	0.29	0
2	FAD	A	501	-	58,58,58	0.30	0	85,89,89	0.30	0
4	GOL	A	503	-	5,5,5	0.93	0	5,5,5	1.07	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	YK6	A	502	-	-	13/17/33/33	0/2/2/2
2	FAD	B	700	-	-	4/34/50/50	0/6/6/6
2	FAD	A	501	-	-	4/34/50/50	0/6/6/6
4	GOL	A	503	-	-	0/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	YK6	O21-C10	9.96	1.60	1.46
3	A	502	YK6	O20-C14	8.92	1.60	1.46
3	A	502	YK6	C10-C09	-2.25	1.44	1.47
3	A	502	YK6	O20-C13	-2.23	1.41	1.45
3	A	502	YK6	O21-C09	-2.19	1.41	1.45

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	YK6	O21-C09-C10	13.68	68.52	59.88
3	A	502	YK6	O20-C13-C14	11.28	68.51	59.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	YK6	O21-C10-C09	-5.84	54.84	59.25
3	A	502	YK6	C10-O21-C09	-5.61	56.64	60.87
3	A	502	YK6	C14-O20-C13	-4.30	56.76	60.78
3	A	502	YK6	O20-C13-C12	-4.17	110.61	117.85
3	A	502	YK6	O20-C14-C13	-3.88	54.73	58.95
3	A	502	YK6	O21-C09-C08	-3.03	112.59	117.85
3	A	502	YK6	C08-C09-C10	-2.48	121.74	125.34

There are no chirality outliers.

All (21) torsion outliers are listed below:

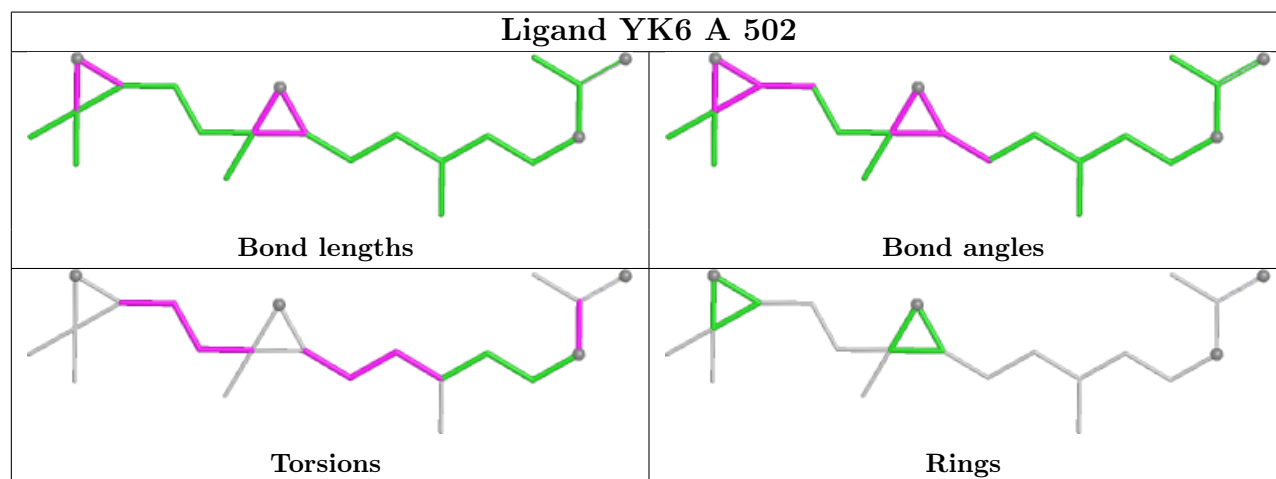
Mol	Chain	Res	Type	Atoms
2	A	501	FAD	C5'-O5'-P-O2P
2	A	501	FAD	C5'-O5'-P-O3P
2	B	700	FAD	C5'-O5'-P-O2P
2	B	700	FAD	C5'-O5'-P-O3P
3	A	502	YK6	C01-C02-O03-C04
3	A	502	YK6	C06-C07-C08-C09
3	A	502	YK6	O16-C02-O03-C04
3	A	502	YK6	C10-C11-C12-C13
3	A	502	YK6	C17-C06-C07-C08
2	A	501	FAD	PA-O3P-P-O5'
2	B	700	FAD	PA-O3P-P-O5'
3	A	502	YK6	C05-C06-C07-C08
3	A	502	YK6	C07-C08-C09-C10
3	A	502	YK6	C18-C10-C11-C12
3	A	502	YK6	O21-C10-C11-C12
3	A	502	YK6	C11-C12-C13-C14
3	A	502	YK6	C09-C10-C11-C12
2	A	501	FAD	P-O3P-PA-O2A
2	B	700	FAD	P-O3P-PA-O2A
3	A	502	YK6	C07-C08-C09-O21
3	A	502	YK6	C11-C12-C13-O20

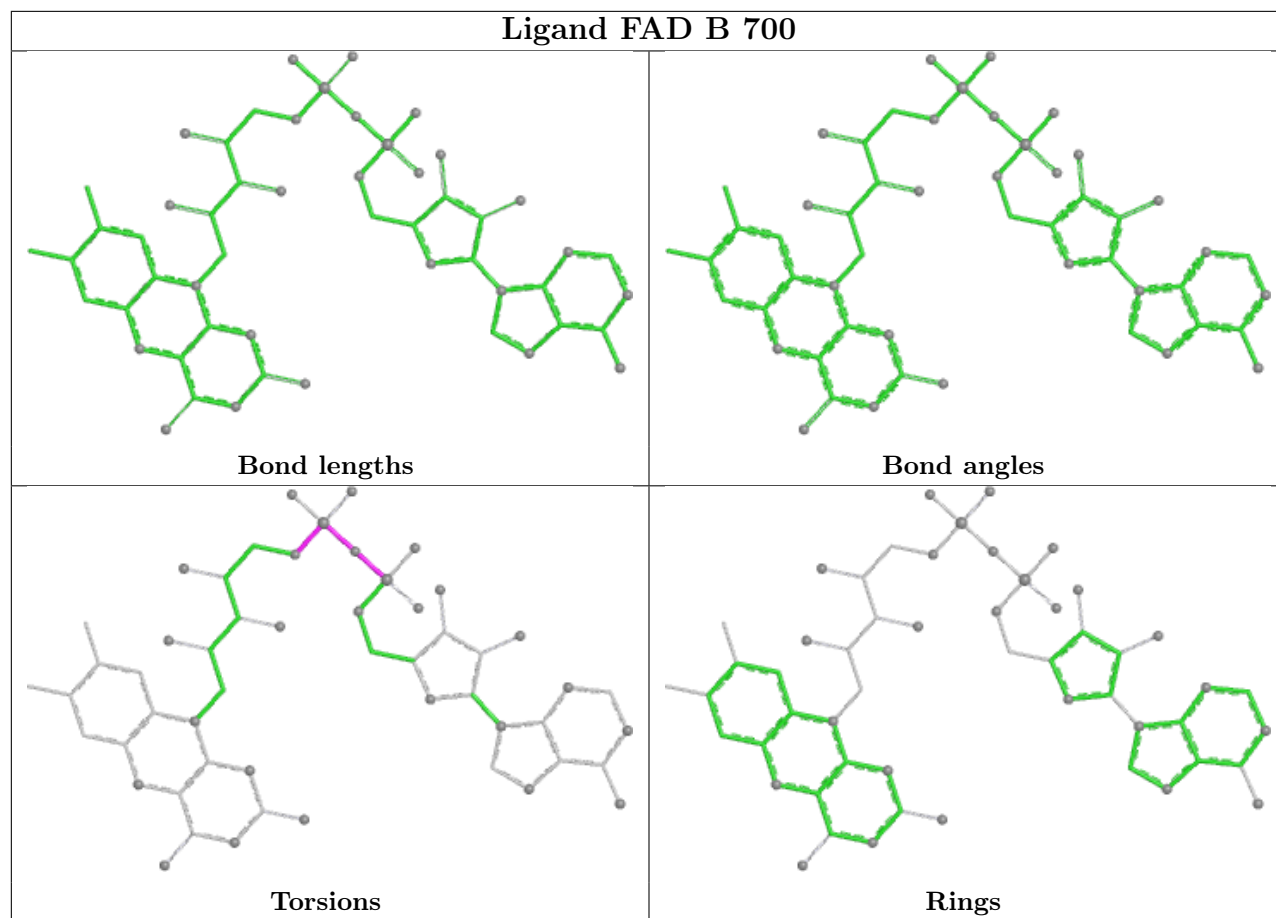
There are no ring outliers.

2 monomers are involved in 3 short contacts:

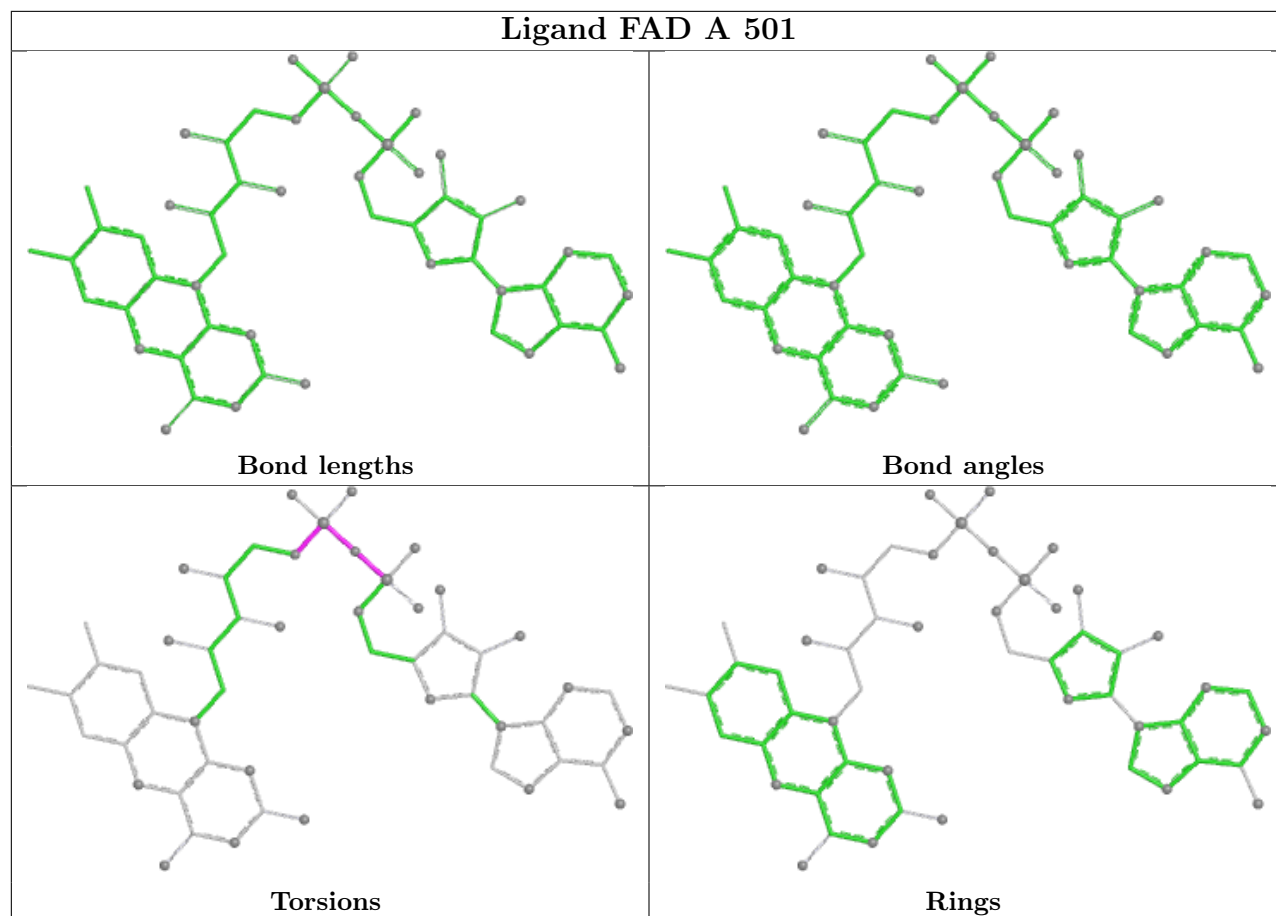
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	700	FAD	1	0
2	A	501	FAD	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

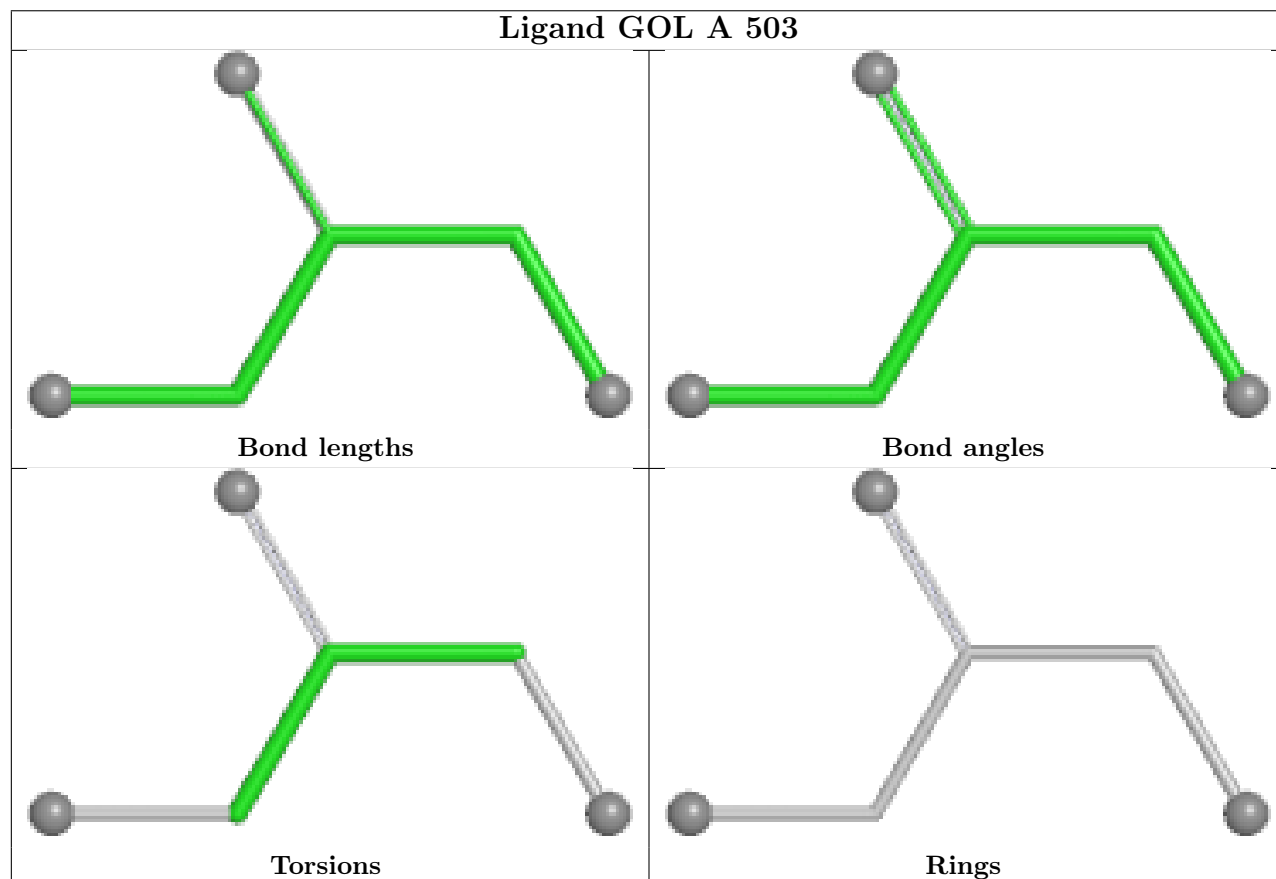




Ligand FAD A 501



Ligand GOL A 503



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	472/511 (92%)	0.02	13 (2%)	55 48	20, 39, 67, 97	1 (0%)
1	B	472/511 (92%)	0.65	35 (7%)	20 15	35, 63, 96, 116	0
All	All	944/1022 (92%)	0.34	48 (5%)	33 26	20, 49, 90, 116	1 (0%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	480	ALA	7.8
1	A	480	ALA	5.6
1	B	383	ALA	5.6
1	B	382	LEU	4.7
1	B	231	ASP	4.6
1	B	208	TYR	4.4
1	A	479	ALA	4.2
1	A	355	GLN	3.9
1	B	363	ALA	3.8
1	B	479	ALA	3.7
1	B	151	GLY	3.4
1	A	362	ARG	3.3
1	B	361	ALA	3.2
1	B	230	ALA	3.1
1	B	362	ARG	3.1
1	A	361	ALA	3.0
1	B	354	VAL	3.0
1	B	4	THR	3.0
1	B	150	SER	2.9
1	B	451	MET	2.9
1	A	363	ALA	2.9
1	B	158	VAL	2.8
1	B	168	GLU	2.7
1	B	166	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	147	GLY	2.6
1	B	355	GLN	2.6
1	B	143	LEU	2.5
1	A	164	SER	2.5
1	A	478	ASP	2.5
1	A	149	GLY	2.5
1	A	4	THR	2.4
1	B	35	VAL	2.4
1	B	317	GLY	2.4
1	B	312	GLU	2.4
1	A	364	GLY	2.4
1	B	229	ALA	2.3
1	B	5	ARG	2.3
1	B	165	GLY	2.2
1	A	208	TYR	2.2
1	B	152	GLY	2.2
1	A	150	SER	2.1
1	B	477	LEU	2.1
1	B	157	VAL	2.1
1	B	478	ASP	2.1
1	B	365	THR	2.1
1	B	291	PRO	2.1
1	B	282	LEU	2.0
1	B	404	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

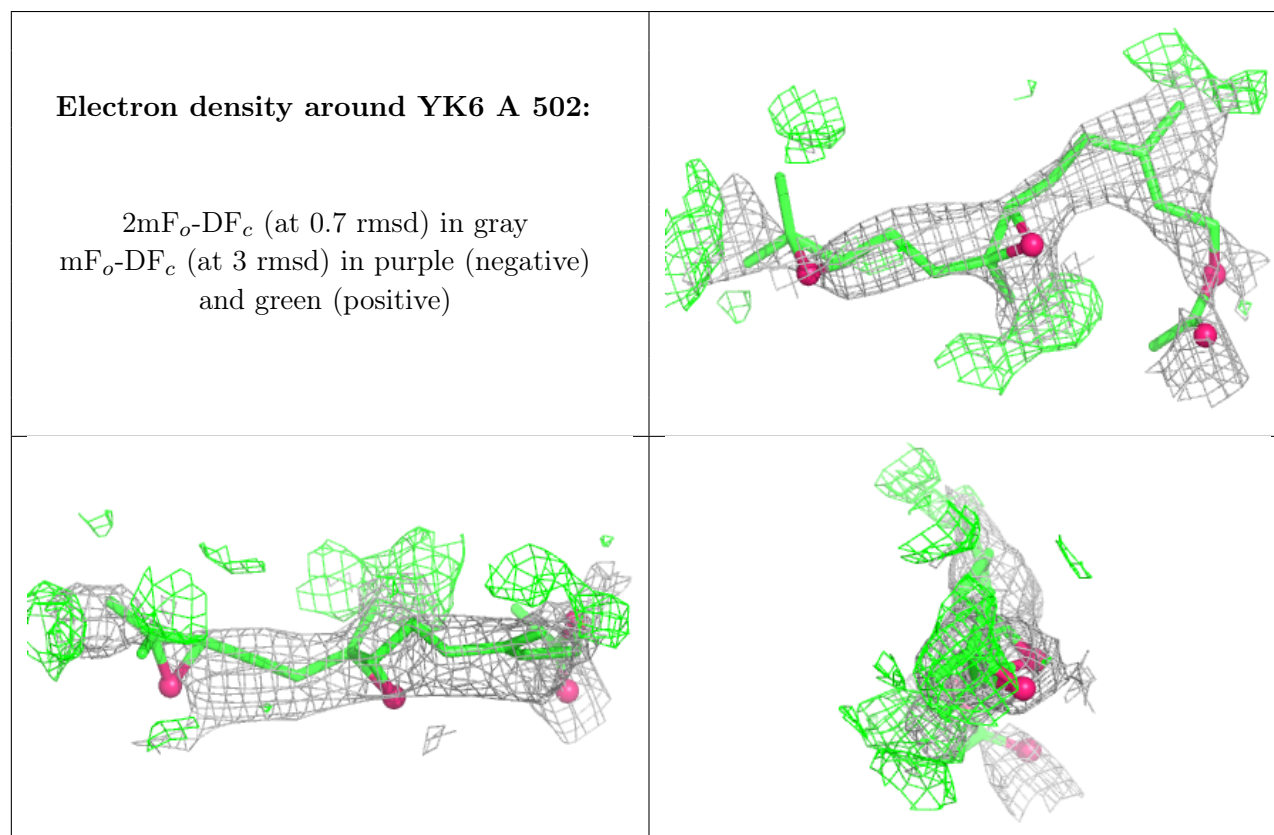
There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

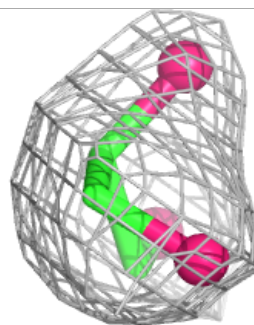
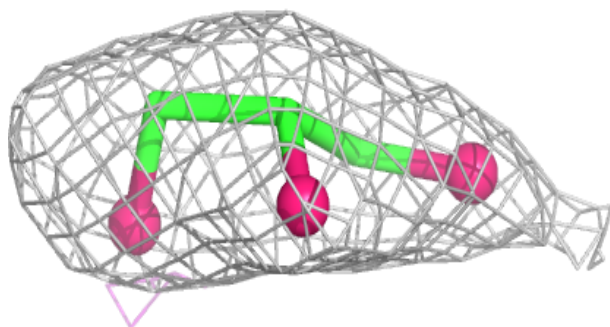
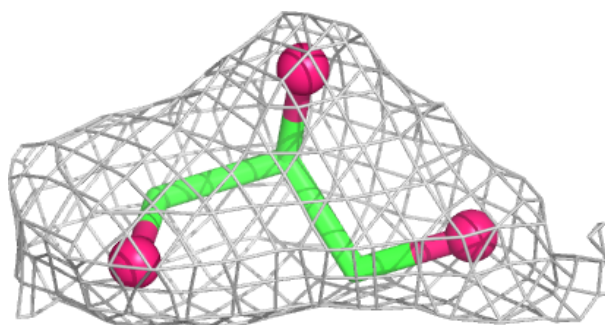
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	YK6	A	502	21/21	0.67	0.29	58,83,93,99	0
4	GOL	A	503	6/6	0.87	0.13	55,58,69,75	0
2	FAD	B	700	53/53	0.95	0.08	41,53,71,81	0
2	FAD	A	501	53/53	0.96	0.08	21,31,48,58	0
5	CL	A	504	1/1	0.97	0.10	42,42,42,42	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



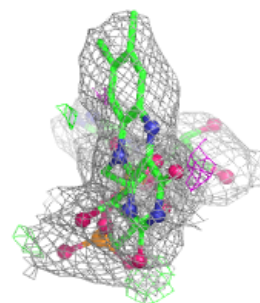
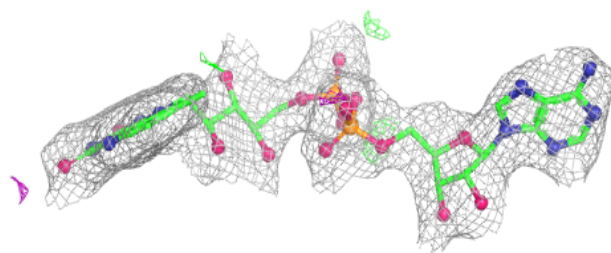
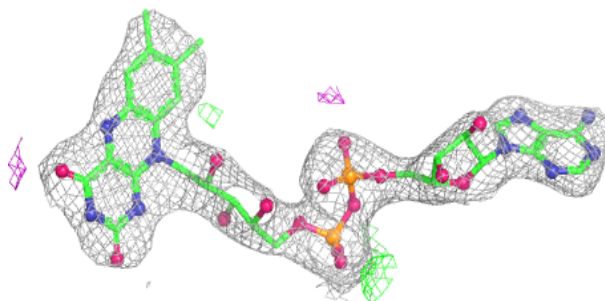
Electron density around GOL A 503:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

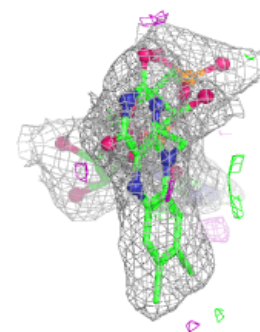
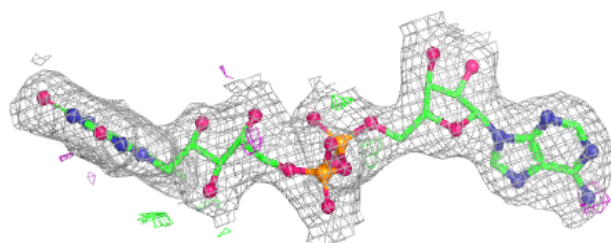
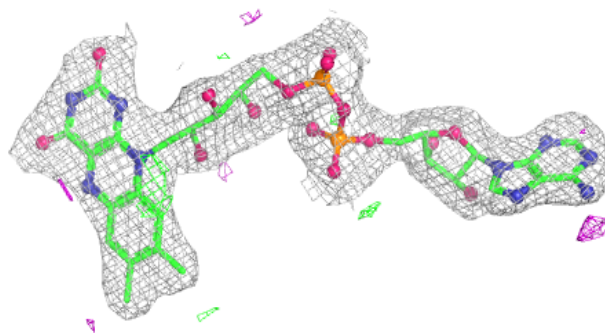


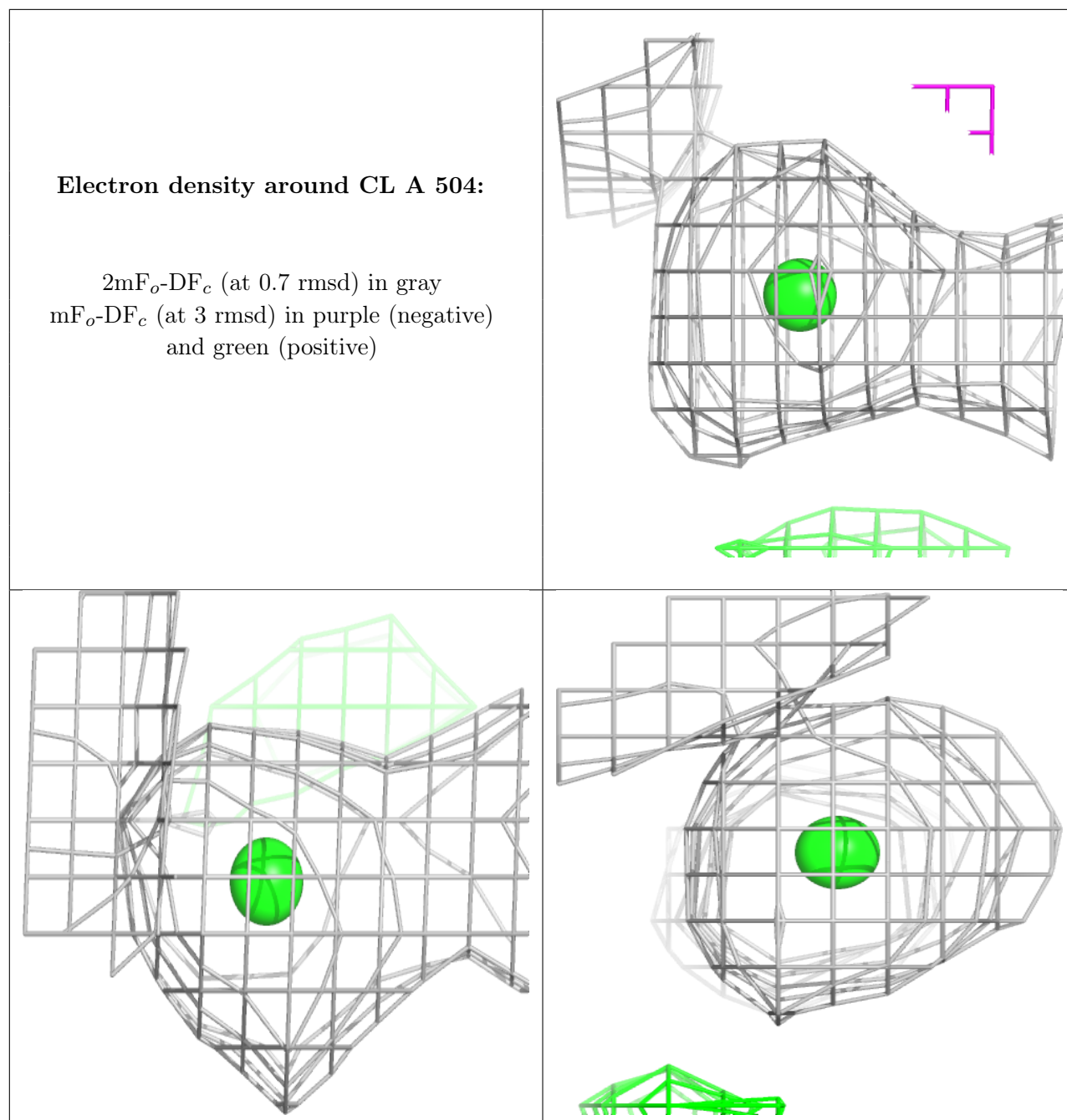
Electron density around FAD B 700:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around FAD A 501:**

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers ⓘ

There are no such residues in this entry.